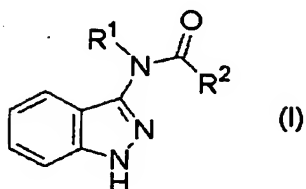


Claims

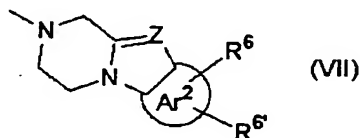
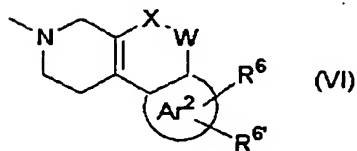
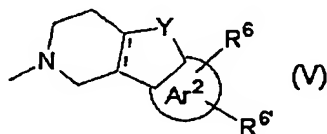
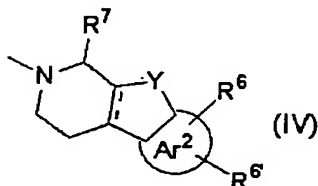
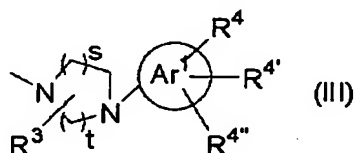
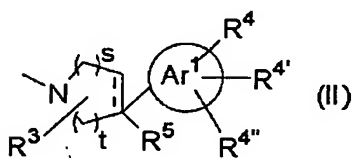
1. An indazole compound represented by the following formula (I):



wherein

R¹ is a hydrogen atom, an optionally substituted alkyl, an optionally substituted phenyl or an optionally substituted aromatic heterocyclic ring, and

R² is any of the following formula (II) to the following formula (VII),



wherein

in the formula (II),

is a single bond or a double bond,

in the formulas (II) and (III),

s is an integer of 1 or 2,

t is an integer of 1 or 2,

R^3 is a hydrogen atom, a halogen atom, an optionally substituted alkyl, a hydroxyl, an alkoxy, a carboxy or an alkoxycarbonyl,
 ring Ar^1 is an aryl or an aromatic heterocyclic ring,
 R^4 , $R^{4'}$, $R^{4''}$ are the same or different and each is a hydrogen atom, a halogen atom, an optionally substituted alkyl, an optionally substituted alkenyl, an optionally substituted alkynyl, a hydroxyl, an alkoxy, a carboxy, an alkoxycarbonyl, an acyl, $-O(C=O)R^{4a}$ (wherein R^{4a} is an optionally substituted C_{1-6} alkyl), $-(C=O)NR^{4a'}R^{4a''}$ (wherein $R^{4a'}$ and $R^{4a''}$ are the same or different and each is a hydrogen atom or an optionally substituted C_{1-6} alkyl, or $R^{4a'}$ and $R^{4a''}$ are taken together to form an optionally substituted 5- to 7-membered non-aromatic heterocyclic ring), $-NH(C=O)R^{4a}$ (wherein R^{4a} is an optionally substituted C_{1-6} alkyl), $-SO_2NR^{4a'}R^{4a''}$ (wherein $R^{4a'}$ and $R^{4a''}$ are the same or different and each is a hydrogen atom or an optionally substituted C_{1-6} alkyl, or $R^{4a'}$ and $R^{4a''}$ are taken together to form an optionally substituted 5- to 7-membered non-aromatic heterocyclic ring), $-NHSO_2R^{4a}$ (wherein R^{4a} is an optionally substituted C_{1-6} alkyl), an amino, an alkylamino, $-SR^{4a}$ (wherein R^{4a} is an optionally substituted C_{1-6} alkyl), $-SO_2R^{4a}$ (wherein R^{4a} is an optionally substituted C_{1-6} alkyl), a cyano, an optionally substituted phenyl or an optionally substituted heterocyclic ring, or R^4 and $R^{4'}$ are taken together to form an C_{1-3} alkylenedioxy, and
 R^5 is absent, or a hydrogen atom, a halogen atom, an optionally substituted alkyl, a hydroxyl, an alkoxy, an alkoxycarbonyl, an acyl, $-(C=O)NR^{5a}R^{5a'}$ (wherein R^{5a} and $R^{5a'}$ are the same or different and each is a hydrogen atom or an optionally substituted C_{1-6} alkyl), $-NH(C=O)R^{5a''}$ (wherein $R^{5a''}$ is an optionally substituted C_{1-6} alkyl), an amino, an alkylamino, $-SR^{5a}$ (wherein R^{5a} is a hydrogen atom or an optionally substituted C_{1-6} alkyl) or a cyano,

in the formulas (IV) and (V),

is a single bond or a double bond,

Y is a carbonyl, NR^{10} , an oxygen atom or a sulfur atom,
5 wherein R^{10} is a hydrogen atom, an optionally substituted
alkyl, an acyl, an alkoxycarbonyl or $-\text{SO}_2\text{R}^{10a}$ (wherein R^{10a}
is an optionally substituted C_{1-6} alkyl or an optionally
substituted phenyl),

ring Ar^2 is a phenyl or an aromatic heterocyclic ring,

10 R^6 and $\text{R}^{6'}$ are the same or different and each is a hydrogen
atom, a halogen atom, an optionally substituted alkyl, an
optionally substituted alkenyl, an optionally substituted
alkynyl, a hydroxyl, an alkoxy, a carboxy, an

alkoxycarbonyl, an acyl, $-\text{O}(\text{C}=\text{O})\text{R}^{6a}$ (wherein R^{6a} is an
15 optionally substituted C_{1-6} alkyl), $-(\text{C}=\text{O})\text{NR}^{6a'}\text{R}^{6a''}$ (wherein
 $\text{R}^{6a'}$ and $\text{R}^{6a''}$ are the same or different and each is a

hydrogen atom or an optionally substituted C_{1-6} alkyl, or
 $\text{R}^{6a'}$ and $\text{R}^{6a''}$ are taken together to form an optionally

substituted 5- to 7-membered non-aromatic heterocyclic

20 ring), $-\text{NH}(\text{C}=\text{O})\text{R}^{6a}$ (wherein R^{6a} is an optionally substituted
 C_{1-6} alkyl), $-\text{SO}_2\text{NR}^{6a'}\text{R}^{6a''}$ (wherein $\text{R}^{6a'}$ and $\text{R}^{6a''}$ are the same
or different and each is a hydrogen atom or an optionally
substituted C_{1-6} alkyl, or $\text{R}^{6a'}$ and $\text{R}^{6a''}$ are taken together to

25 form an optionally substituted 5- to 7-membered non-
aromatic heterocyclic ring), $-\text{NHSO}_2\text{R}^{6a}$ (wherein R^{6a} is an
optionally substituted C_{1-6} alkyl), an amino, an

alkylamino, $-\text{SR}^{6a}$ (wherein R^{6a} is an optionally substituted
 C_{1-6} alkyl), a cyano, an optionally substituted phenyl or
an optionally substituted heterocyclic ring, or

30 R^4 and $\text{R}^{4'}$ are taken together to form a C_{1-3} alkylenedioxy,
and

R^7 is a hydrogen atom or an optionally substituted alkyl,
in the formula (VI),

35 X and W are any of $\text{C}(=\text{O})$ and O, $\text{C}(=\text{O})$ and NR^{11} , and NR^{11} and
 $\text{C}(=\text{O})$,

wherein R^{11} is a hydrogen atom or an optionally substituted alkyl,

ring Ar^2 is a phenyl or an aromatic heterocyclic ring, and R^6 and $R^{6'}$ are the same or different and each is a hydrogen

5 atom, a halogen atom, an optionally substituted alkyl, an optionally substituted alkenyl, an optionally substituted alkynyl, a hydroxyl, an alkoxy, a carboxy, an

alkoxycarbonyl, an acyl, $-O(C=O)R^{6a}$ (wherein R^{6a} is an optionally substituted C_{1-6} alkyl), $-(C=O)NR^{6a'}R^{6a''}$ (wherein $R^{6a'}$ and $R^{6a''}$ are the same or different and each is a hydrogen atom or an optionally substituted C_{1-6} alkyl, or $R^{6a'}$ and $R^{6a''}$ are taken together to form an optionally substituted 5- to 7-membered non-aromatic heterocyclic

ring), $-NH(C=O)R^{6a}$ (wherein R^{6a} is an optionally substituted C_{1-6} alkyl), $-SO_2NR^{6a'}R^{6a''}$ (wherein $R^{6a'}$ and $R^{6a''}$ are the same or different and each is a hydrogen atom or an optionally substituted C_{1-6} alkyl, or $R^{6a'}$ and $R^{6a''}$ are taken together to form an optionally substituted 5- to 7-membered non-

aromatic heterocyclic ring), $-NHSO_2R^{6a}$ (wherein R^{6a} is an optionally substituted C_{1-6} alkyl), an amino, an alkylamino, $-SR^{6a}$ (wherein R^{6a} is an optionally substituted C_{1-6} alkyl), a cyano, an optionally substituted phenyl or an optionally substituted heterocyclic ring, or

R^4 and $R^{4'}$ are taken together to form a C_{1-3} alkylenedioxy, and

in the formula (VII),

Z is a carbon atom or a nitrogen atom,

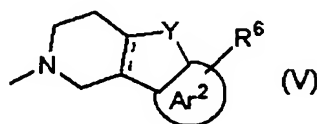
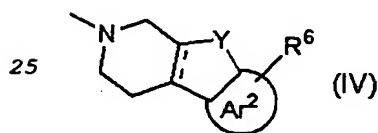
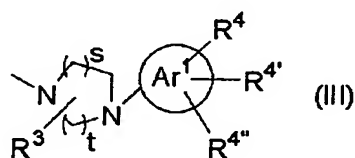
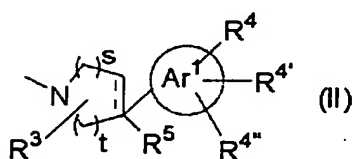
ring Ar^2 is a phenyl or an aromatic heterocyclic ring, and R^6 and $R^{6'}$ are the same or different and each is a hydrogen

30 atom, a halogen atom, an optionally substituted alkyl, an optionally substituted alkenyl, an optionally substituted alkynyl, a hydroxyl, an alkoxy, a carboxy, an

alkoxycarbonyl, an acyl, $-O(C=O)R^{6a}$ (wherein R^{6a} is an optionally substituted C_{1-6} alkyl), $-(C=O)NR^{6a'}R^{6a''}$ (wherein $R^{6a'}$ and $R^{6a''}$ are the same or different and each is a

hydrogen atom or an optionally substituted C₁₋₆ alkyl, or R^{6a'} and R^{6a''} are taken together to form an optionally substituted 5- to 7-membered non-aromatic heterocyclic ring), -NH(C=O)R^{6a} (wherein R^{6a} is an optionally substituted C₁₋₆ alkyl), -SO₂NR^{6a'}R^{6a''} (wherein R^{6a'} and R^{6a''} are the same or different and each is a hydrogen atom or an optionally substituted C₁₋₆ alkyl, or R^{6a'} and R^{6a''} are taken together to form an optionally substituted 5- to 7-membered non-aromatic heterocyclic ring), -NHSO₂R^{6a} (wherein R^{6a} is an optionally substituted C₁₋₆ alkyl), an amino, an alkylamino, -SR^{6a} (wherein R^{6a} is an optionally substituted C₁₋₆ alkyl), a cyano, an optionally substituted phenyl or an optionally substituted heterocyclic ring, or R⁴ and R^{4'} are taken together to form a C₁₋₃ alkylenedioxy, a pharmaceutically acceptable salt thereof, a hydrate thereof, a water adduct thereof or a solvate thereof.

2. The indazole compound of claim 1, wherein, in the above-mentioned formula (I), R² is any of the following formula (II) to the following formula (V),



wherein
in the formula (II),

is a single bond or a double bond,
in the formulas (II) and (III),

s is an integer of 1 or 2,
t is an integer of 0 to 2,
R³ is a hydrogen atom, a halogen atom, an optionally
substituted alkyl, a carboxyl, an alkoxycarbonyl, a
5 hydroxy or an alkoxy,
ring Ar¹ is a phenyl or an aromatic heterocyclic ring,
R⁴, R^{4'} and R^{4''} are the same or different and each is a
hydrogen atom, a halogen atom, an optionally substituted
alkyl, an alkoxycarbonyl, a hydroxy, an alkoxy, a
10 sulfonamide, a mercapto, a sulfinyl, a sulfonyl, an amino
or an alkylamino, and
R⁵ is absent, or a hydrogen atom, a halogen atom, an
optionally substituted alkyl, a hydroxy, an alkoxy, an
amino, an alkylamino, a sulfanyl or a cyano, and
15 in the formulas (IV) and (V),

is a single bond or a double bond,
Y is a carbonyl, NR¹⁰, an oxygen atom or a sulfur atom,
20 wherein R¹⁰ is a hydrogen atom, an optionally substituted
alkyl, an acyl, an alkoxycarbonyl or a sulfonyl,
ring Ar² is a phenyl or an aromatic heterocyclic ring,
R⁶ is a hydrogen atom, a halogen atom, an optionally
substituted alkyl, a cyano, a hydroxy or an alkoxy,
25 a pharmaceutically acceptable salt thereof, a hydrate thereof,
a water adduct thereof or a solvate thereof.

3. The indazole compound of claim 1,
wherein,
30 in the above-mentioned formula (I),
R¹ is a hydrogen atom or an optionally substituted alkyl,
in the above-mentioned formulas (II) and (III),
s is an integer of 1,
t is an integer of 2,
35 R³ is a hydrogen atom,

ring Ar¹ is a phenyl or a thiophene,
R⁴, R^{4'}, R^{4''} are the same or different and each is a hydrogen
atom, a halogen atom, an optionally substituted alkyl, a
hydroxy, an alkoxy, -SR^{4a} (wherein R^{4a} is an optionally
5 substituted C₁₋₆ alkyl) or a cyano, and
R⁵ is a hydroxy or a cyano,
in the above-mentioned formulas (IV) and (V),
Y is NR¹⁰,

wherein R¹⁰ is a hydrogen atom or an optionally substituted
10 alkyl,

ring Ar² is a phenyl, and
R⁶ and R^{6'} are the same or different and each is a hydrogen
atom, a halogen atom, an optionally substituted alkyl, a
hydroxy or an alkoxy,
15 in the above-mentioned formula (VI),
X and W are any of C(=O) and O, C(=O) and NR¹¹, and NR¹¹ and
C(=O),

wherein R¹¹ is a hydrogen atom,
ring Ar² is a phenyl, and
20 R⁶ and R^{6'} are the same or different and each is a hydrogen
atom, a halogen atom or an optionally substituted alkyl, and
in the above-mentioned formula (VII),
ring Ar² is a phenyl, and
R⁶ and R^{6'} are the same or different and each is a hydrogen
25 atom, a halogen atom or an optionally substituted alkyl,
a pharmaceutically acceptable salt thereof, a hydrate thereof,
a water adduct thereof or a solvate thereof.

4. The indazole compound of claim 1 or 3,
30 wherein,
in the above-mentioned formula (I),
R¹ is a hydrogen atom,
in the above-mentioned formulas (II) and (III),
s is an integer of 1,
35 t is an integer of 2,

R³ is a hydrogen atom,
ring Ar¹ is a phenyl,
R⁴, R^{4'}, R^{4''} are the same or different and each is a hydrogen
atom, a halogen atom or an optionally substituted alkyl, and
5 R⁵ is a hydroxy or a cyano, and
in the above-mentioned formula (IV),
Y is NR¹⁰,

wherein R¹⁰ is a hydrogen atom or a methyl,
a pharmaceutically acceptable salt thereof, a hydrate thereof,
10 a water adduct thereof or a solvate thereof.

5. The indazole compound of any of claims 1 to 4,
wherein,
in the above-mentioned formula (I),
15 R¹ is a hydrogen atom, and
in the above-mentioned formula (II),
s is an integer of 1,
t is an integer of 2,
R³ is a hydrogen atom,
20 ring Ar¹ is a phenyl,
R⁴, R^{4'}, R^{4''} are the same or different and each is a hydrogen
atom, a halogen atom or an optionally substituted alkyl, and
R⁵ is a hydroxyl,
a pharmaceutically acceptable salt thereof, a hydrate thereof,
25 a water adduct thereof or a solvate thereof.

6. The indazole compound of claim 1, which is selected from
(1) 4-[4-chloro-3-(trifluoromethyl)phenyl]-4-hydroxy-1-
piperidinecarboxylic acid (1H-indazol-3-yl)amide,
30 (3) 4-hydroxy-4-[3-(trifluoromethyl)phenyl]-1-
piperidinecarboxylic acid (1H-indazol-3-yl)amide,
(4) 4-(4-chlorophenyl)-4-hydroxy-1-piperidinecarboxylic acid
(1H-indazol-3-yl)amide,
(6) 4-[3-fluoro-5-(trifluoromethyl)phenyl]-4-hydroxy-1-
35 piperidinecarboxylic acid (1H-indazol-3-yl)amide,

- (9) 4-[4-fluoro-3-(trifluoromethyl)phenyl]-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
- (10) 4-hydroxy-4-[4-methyl-3-(trifluoromethyl)phenyl]-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
- 5 (12) 4-(3,5-difluorophenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
- (15) 4-(3-chloro-4-fluorophenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
- (20) 4-(3-chloro-2-fluorophenyl)-4-hydroxy-1-
- 10 piperidinecarboxylic acid (1H-indazol-3-yl)amide,
- (21) 4-(3,4-dichlorophenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
- (22) 4-(3-chloro-5-fluorophenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
- 15 (23) 4-(4-chloro-3-methylphenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
- (24) 4-(3-chlorophenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
- (27) 4-(1,3-benzodioxol-5-yl)-4-hydroxy-1-piperidinecarboxylic
- 20 acid (1H-indazol-3-yl)amide,
- (28) 4-hydroxy-4-(3-methylphenyl)-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
- (29) 4-(3-cyanophenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
- 25 (30) 4-hydroxy-4-[3-(methylthio)phenyl]-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
- (31) 4-(3-ethylphenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
- (33) 4-(2,5-dichlorophenyl)-4-hydroxy-1-piperidinecarboxylic
- 30 acid (1H-indazol-3-yl)amide,
- (34) 4-[3,5-bis(trifluoromethyl)phenyl]-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
- (35) 4-[2-fluoro-5-(trifluoromethyl)phenyl]-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
- 35 (36) 4-[2-chloro-5-(trifluoromethyl)phenyl]-4-hydroxy-1-

- piperidinecarboxylic acid (1H-indazol-3-yl) amide,
- (40) 4-cyano-4-(2-methoxyphenyl)-1-piperidinecarboxylic acid
(1H-indazol-3-yl) amide,
- (42) 4-cyano-4-[3-(trifluoromethyl)phenyl]-1-
- 5 piperidinecarboxylic acid (1H-indazol-3-yl) amide,
- (43) 4-cyano-4-(2-fluorophenyl)-1-piperidinecarboxylic acid
(1H-indazol-3-yl) amide,
- (44) 4-[4-chloro-3-(trifluoromethyl)phenyl]-4-cyano-1-
- piperidinecarboxylic acid (1H-indazol-3-yl) amide,
- 10 (46) 4-(5-bromo-2-thienyl)-4-cyano-1-piperidinecarboxylic acid
(1H-indazol-3-yl) amide,
- (47) 4-cyano-4-(3,5-difluorophenyl)-1-piperidinecarboxylic
acid (1H-indazol-3-yl) amide,
- (48) 4-(4-bromo-2-chlorophenyl)-4-cyano-1-piperidinecarboxylic
- 15 acid (1H-indazol-3-yl) amide,
- (49) 4-phenyl-1,2,3,6-tetrahydropyridine-1-carboxylic acid
(1H-indazol-3-yl) amide,
- (50) 4-(4-fluorophenyl)-1,2,3,6-tetrahydropyridine-1-
- carboxylic acid (1H-indazol-3-yl) amide,
- 20 (52) 4-(2-fluorophenyl)-1,2,3,6-tetrahydropyridine-1-
- carboxylic acid (1H-indazol-3-yl) amide,
- (53) 4-(3-chloro-4-fluorophenyl)-1,2,3,6-tetrahydropyridine-1-
- carboxylic acid (1H-indazol-3-yl) amide,
- (55) 4-(3-fluorophenyl)-1,2,3,6-tetrahydropyridine-1-
- 25 carboxylic acid (1H-indazol-3-yl) amide,
- (56) 4-(2,3-difluorophenyl)-1,2,3,6-tetrahydropyridine-1-
- carboxylic acid (1H-indazol-3-yl) amide,
- (58) 4-(5-chloro-2-thienyl)-1,2,3,6-tetrahydropyridine-1-
- carboxylic acid (1H-indazol-3-yl) amide,
- 30 (59) 4-(3-methyl-2-thienyl)-1,2,3,6-tetrahydropyridine-1-
- carboxylic acid (1H-indazol-3-yl) amide,
- (60) 4-(2-thienyl)-1,2,3,6-tetrahydropyridine-1-carboxylic
acid (1H-indazol-3-yl) amide,
- (61) 4-[3-(trifluoromethyl)phenyl]-1,2,3,6-tetrahydropyridine-
- 35 1-carboxylic acid (1H-indazol-3-yl) amide,

- (62) 4-(3,4-dimethoxyphenyl)-1,2,3,6-tetrahydropyridine-1-carboxylic acid (1H-indazol-3-yl)amide,
- (63) 4-[3-(dimethylamino)phenyl]-1,2,3,6-tetrahydropyridine-1-carboxylic acid (1H-indazol-3-yl)amide,
- 5 (64) 1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid (1H-indazol-3-yl)amide,
- (65) 9-methyl-1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid (1H-indazol-3-yl)amide,
- (66) 9-(2-methoxyethyl)-1,3,4,9-tetrahydro- β -carboline-2-
- 10 carboxylic acid (1H-indazol-3-yl)amide,
- (69) 6-(trifluoromethyl)-1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid (1H-indazol-3-yl)amide,
- (70) 6-fluoro-1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid (1H-indazol-3-yl)amide,
- 15 (71) 7-fluoro-1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid (1H-indazol-3-yl)amide,
- (72) 6-chloro-1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid (1H-indazol-3-yl)amide,
- (73) 6-methoxy-1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid
- 20 (1H-indazol-3-yl)amide,
- (74) 6-hydroxy-1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid (1H-indazol-3-yl)amide,
- (75) 7-chloro-1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid (1H-indazol-3-yl)amide,
- 25 (76) 7-(trifluoromethyl)-1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid (1H-indazol-3-yl)amide,
- (77) 5-fluoro-1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid (1H-indazol-3-yl)amide,
- (78) 5-chloro-1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid
- 30 (1H-indazol-3-yl)amide,
- (79) 8-methyl-1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid (1H-indazol-3-yl)amide,
- (80) 3,4-dihydro[1]benzothieno[2,3-c]pyridine-2-carboxylic acid (1H-indazol-3-yl)amide,
- 35 (81) 6-methyl-1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid

- (1H-indazol-3-yl) amide,
- (82) 7-chloro-6-fluoro-1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid (1H-indazol-3-yl) amide,
- (83) 7-chloro-6-(trifluoromethyl)-1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid (1H-indazol-3-yl) amide,
- 5 (93) 4-[4-chloro-3-(trifluoromethyl)phenyl]-1-piperazinecarboxylic acid (1H-indazol-3-yl) amide,
- (94) 4-[4-fluoro-3-(trifluoromethyl)phenyl]-1-piperazinecarboxylic acid (1H-indazol-3-yl) amide,
- 10 (95) 4-[4-methoxy-3-(trifluoromethyl)phenyl]-1-piperazinecarboxylic acid (1H-indazol-3-yl) amide,
- (97) 4-[3-fluoro-5-(trifluoromethyl)phenyl]-1-piperazinecarboxylic acid (1H-indazol-3-yl) amide,
- (98) 4-(3,4-dichlorophenyl)-1-piperazinecarboxylic acid (1H-indazol-3-yl) amide,
- 15 (99) 4-[2-chloro-5-(trifluoromethyl)phenyl]-1-piperazinecarboxylic acid (1H-indazol-3-yl) amide,
- (100) 4-[3-(trifluoromethyl)phenyl]-1-piperazinecarboxylic acid (1H-indazol-3-yl) amide,
- 20 (103) 5-oxo-1,5-dihydro-2H-chromeno[3,4-c]pyridine-3-carboxylic acid (1H-indazol-3-yl) amide,
- (104) 5-oxo-1,4,5,6-tetrahydrobenzo[c]-2,7-naphthyridine-3-carboxylic acid (1H-indazol-3-yl) amide,
- (105) 3,4-dihydropyrazino[1,2-a]benzimidazole-2-carboxylic acid (1H-indazol-3-yl) amide,
- 25 (106) 3,4-dihydropyrazino[1,2-a]indole-2-carboxylic acid (1H-indazol-3-yl) amide,
- (108) 1-[(dimethylamino)methyl]-1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid (1H-indazol-3-yl) amide,
- 30 (109) 6-oxo-1,4,5,6-tetrahydrobenzo[c]-1,7-naphthyridine-3-carboxylic acid (1H-indazol-3-yl) amide,
- (112) 4-[3-(trifluoromethyl)phenyl]piperidine-1-carboxylic acid (1H-indazol-3-yl) amide,
- (116) 4-[4-chloro-3-(trifluoromethyl)phenyl]-4-methoxypiperidine-1-carboxylic acid (1H-indazol-3-yl) amide,
- 35

(117) 4-[4-chloro-3-(trifluoromethyl)phenyl]-3-methylpiperazine-1-carboxylic acid (1H-indazol-3-yl)amide,
(123) 4-[4-chloro-3-(trifluoromethyl)phenyl]-4-fluoropiperidine-1-carboxylic acid (1H-indazol-3-yl)amide,
5 (130) 4-(2-fluoro-5-methylphenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
(131) 4-(3-chloro-2-methylphenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
(132) 4-(3-chloro-4-methylphenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
10 (134) 4-(3-fluoro-2-methylphenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
(135) 4-(5-fluoro-2-methylphenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
15 (136) 4-(4-fluoro-3-methylphenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
(138) 4-(3-fluoro-5-methylphenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
(139) 4-(2,5-dimethylphenyl)-4-hydroxy-1-piperidinecarboxylic
20 acid (1H-indazol-3-yl)amide,
(140) 4-hydroxy-4-[2-methyl-3-(trifluoromethyl)phenyl]-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
(141) 4-hydroxy-4-[2-methyl-5-(trifluoromethyl)phenyl]-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
25 (142) 4-(3,4-dimethylphenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide,
(143) 4-(3,5-dimethylphenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide, and
(144) 4-(2,3-dimethylphenyl)-4-hydroxy-1-piperidinecarboxylic
30 acid (1H-indazol-3-yl)amide,
a pharmaceutically acceptable salt thereof, a hydrate thereof,
a water adduct thereof or a solvate thereof.

7. The indazole compound of claim 1, which is 4-hydroxy-4-(3-methylphenyl)-1-piperidinecarboxylic acid (1H-indazol-3-

yl)amide, a pharmaceutically acceptable salt thereof, a hydrate thereof, a water adduct thereof or a solvate thereof.

8. The indazole compound of claim 1, which is 4-(3-chloro-2-fluorophenyl)-4-hydroxy-1-piperidinecarboxylic acid (1H-indazol-3-yl)amide, a pharmaceutically acceptable salt thereof, a hydrate thereof, a water adduct thereof or a solvate thereof.

9. The indazole compound of claim 1, which is 4-(4-fluorophenyl)-1,2,3,6-tetrahydropyridine-1-carboxylic acid (1H-indazol-3-yl)amide, a pharmaceutically acceptable salt thereof, a hydrate thereof, a water adduct thereof or a solvate thereof.

10. The indazole compound of claim 1, which is 1,3,4,9-tetrahydro- β -carboline-2-carboxylic acid (1H-indazol-3-yl)amide, a pharmaceutically acceptable salt thereof, a hydrate thereof, a water adduct thereof or a solvate thereof.

11. The indazole compound of claim 1, which is 4-[4-chloro-3-(trifluoromethyl)phenyl]-1-piperazinecarboxylic acid (1H-indazol-3-yl)amide, a pharmaceutically acceptable salt thereof, a hydrate thereof, a water adduct thereof or a solvate thereof.

12. An agent for the prophylaxis and/or treatment of cancer, which comprises an indazole compound of any one of claims 1 to 11, a pharmaceutically acceptable salt thereof, a hydrate thereof, a water adduct thereof or a solvate thereof.